

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101024\
 Data File : BM048008.D
 Acq On : 11 Oct 2024 11:09
 Operator : RC/JU
 Sample : SP6641
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP6641

Quant Time: Oct 11 11:55:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 10 18:01:13 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.796	152	6276	0.400	ng/ul	0.00
4) Naphthalene-d8	10.579	136	19461	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.423	164	10613	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.162	188	18586	0.400	ng/ul	0.00
17) Chrysene-d12	21.330	240	14523	0.400	ng/ul	0.00
23) Perylene-d12	23.598	264	14151	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/ul	
6) 2-Methylnaphthalene-d10	0.000	152	0d	0.000	ng/ul	
18) Fluoranthene-d10	0.000	212	0d	0.000	ng/ul	
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.307	88	39244	4.066	ng/ul#	79
5) Naphthalene	10.628	128	41707	0.786	ng/ul	99
7) 2-Methylnaphthalene	12.251	142	23732	0.746	ng/ul	99
8) 1-Methylnaphthalene	12.465	142	26496	0.794	ng/ul	100
10) Acenaphthylene	14.141	152	35813	0.694	ng/ul	99
11) Acenaphthene	14.483	153	28134	0.794	ng/ul	99
12) Fluorene	15.469	166	28649	0.715	ng/ul	100
14) Pentachlorophenol	16.820	266	9124	1.633	ng/ul	100
15) Phenanthrene	17.204	178	41167	0.746	ng/ul	99
16) Anthracene	17.297	178	31207	0.635	ng/ul	99
19) Fluoranthene	19.216	202	45334	0.742	ng/ul	96
20) Pyrene	19.578	202	49742	0.765	ng/ul	97
21) Benzo(a)anthracene	21.315	228	36924	0.601	ng/ul	99
22) Chrysene	21.365	228	45389	0.711	ng/ul	99
24) Benzo(b)fluoranthene	22.911	252	40946	0.665	ng/ul	94
25) Benzo(k)fluoranthene	22.954	252	48149	0.777	ng/ul#	94
26) Benzo(a)pyrene	23.498	252	38595	0.795	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	25.903	276	51476	0.666	ng/ul#	93
28) Dibenzo(a,h)anthracene	25.913	278	38076	0.653	ng/ul	97
29) Benzo(g,h,i)perylene	26.600	276	44406	0.716	ng/ul	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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